

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE121318\
 Data File : BE098416.D
 Acq On : 14 Dec 2018 2:58
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 14 10:06:54 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE121018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 10 14:59:55 2018
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	72	0.00
2	1,4-Dioxane	0.400	0.391	2.3	72	-0.01
3	n-Nitrosodimethylamine	0.400	0.225	43.8#	44	0.01
4 S	2-Fluorophenol	0.400	0.387	3.3	72	0.00
5 S	Phenol-d6	0.400	0.407	-1.7	73	0.00
6	bis(2-Chloroethyl)ether	0.400	0.366	8.5	68	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	74	-0.01
8 S	Nitrobenzene-d5	0.400	0.349	12.8	71	0.01
9	Naphthalene	0.400	0.384	4.0	73	0.00
10	Hexachlorobutadiene	0.400	0.405	-1.3	77	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.372	7.0	70	0.00
12	2-Methylnaphthalene	0.400	0.365	8.8	69	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	72	0.00
14 S	2,4,6-Tribromophenol	0.400	0.369	7.8	70	0.00
15 S	2-Fluorobiphenyl	0.400	0.364	9.0	72	0.00
16	Acenaphthylene	0.400	0.386	3.5	72	0.00
17	Acenaphthene	0.400	0.378	5.5	69	0.00
18	Fluorene	0.400	0.382	4.5	69	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	71	-0.01
20	4-Bromophenyl-phenylether	0.400	0.348	13.0	64	0.00
21	Hexachlorobenzene	0.400	0.381	4.8	68	0.00
22	Pentachlorophenol	0.400	0.215	46.3#	36	0.01
23	Phenanthrene	0.400	0.381	4.8	68	0.00
24	Anthracene	0.400	0.344	14.0	61	0.00
25 SURR	Fluoranthene-d10	0.400	0.419	-4.7	71	0.00
26	Fluoranthene	0.400	0.427	-6.7	72	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	86	0.00
28	Pvrene	0.400	0.342	14.5	72	0.00
29 S	Terphenyl-d14	0.400	0.326	18.5	71	0.00
30	Benzo(a)anthracene	0.400	0.360	10.0	79	-0.01
31	Chrysene	0.400	0.374	6.5	81	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.320	20.0	69	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.377	5.8	85	0.00
34 I	Perylene-d12	0.400	0.400	0.0	88	-0.01
35	Benzo(b)fluoranthene	0.400	0.397	0.8	90	0.00
36	Benzo(k)fluoranthene	0.400	0.412	-3.0	94	-0.01
37 C	Benzo(a)pyrene	0.400	0.397	0.8	89	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.368	8.0	83	-0.01
39	Benzo(a,h,i)perylene	0.400	0.383	4.3	87	-0.02

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

(#) = Out of Range	SPCC's out = 0 CCC's out = 0				